## Amendments to the Claims

This listing of claims will replace all prior versions, and listings, of claims in the application:

## **Listing of Claims:**

1. (Previously Presented) An indole compound represented by the formula (I), or a pharmaceutically acceptable salt, solvate, or prodrug thereof;

$$R_{5}$$
 $R_{7}$ 
 $R_{1}$ 
 $R_{2}$ 
 $R_{2}$ 

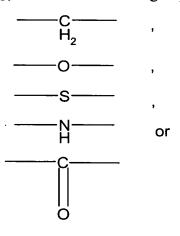
wherein;

R<sub>1</sub> is (c) wherein;

(c) is the group -( $L_1$ )- $R_{11}$ ; where, -( $L_1$ )- is a divalent linking group of 1 to 8 atoms and where  $R_{11}$  is -( $CH_2$ )<sub>m</sub>- $R_{12}$ ;

R2 is hydrogen, or C1-C4 alkyl;

R<sub>3</sub> is -(L<sub>3</sub>)- Z, where -(L<sub>3</sub>)- is a divalent linker group selected from a bond or:



Serial No. 10/629,992

and Z is a group represented by the formulae,

$$R_a$$
 $R_a$ 
 $NH_2$ 

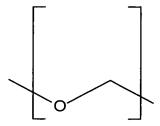
wherein, X is oxygen or sulfur; and R<sub>a</sub> is selected from hydrogen, C<sub>1</sub>-C<sub>8</sub> alkyl, aryl, C<sub>1</sub>-C<sub>8</sub> alkoxy, aralkyl and -CN;

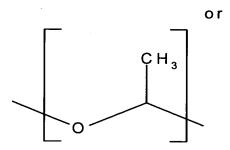
R4 is the group, -(Lh)-(hydroxyfunctional amide); wherein -(Lh)-, is an hydroxyfunctional amide linker having an hydroxyfunctional amide linker length of 1 to 8;

R5 is selected from hydrogen, a non-interfering substituent, or the group,  $-(L_a)$ (acidic group); wherein  $-(L_a)$ -, is an acid linker having an acid linker length of 1 to 8;

 $R_6$  and  $R_7$  are selected from hydrogen,  $C_1$ - $C_6$  alkyl,  $C_2$ - $C_6$  alkenyl, and  $C_2$ - $C_6$  alkynyl.

- 2. (Previously Presented) The compound of claim 1 wherein  $R_2$  is hydrogen,  $C_1$ - $C_4$  alkyl,  $C_2$ - $C_4$  alkenyl, -O-( $C_1$ - $C_3$  alkyl), -S-( $C_1$ - $C_3$  alkyl), and  $C_3$ - $C_4$  cycloalkyl.
  - 3. (Cancelled)
- 4. (Previously Presented) The compound of Claim 1 wherein the hydroxyfunctional amide linker group, -(Lh)-, for R<sub>4</sub> is a divalent group selected from,



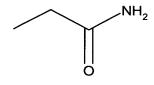


where R<sub>40</sub>, R<sub>41</sub>, R<sub>42</sub>, and R<sub>43</sub> are each independently selected from hydrogen, C<sub>1</sub>-C<sub>8</sub> alkyl.

- 5. (Cancelled)
- 6. (Previously Presented) The compound of claim 1 wherein R<sub>5</sub> is the group, -(L<sub>a</sub>)-(acidic group) and wherein the (acidic group) is:

-COOH.

- 7. (Cancelled)
- 8. (Cancelled)
- 9. (Cancelled)
- 10. (Original) The compound of claim 1 wherein for R<sub>3</sub>, Z is the group represented by the formula;



and the linking group  $-(L_3)$ - is a bond.

- 11. (Cancelled)
- 12. (Cancelled)
- 13. (Cancelled)
- 14. (Cancelled)

- 15. (Cancelled)
- 16. (Cancelled)
- 17. (Cancelled)
- 18. (Original) The compound of claim 1 wherein R4 is the group, -(L<sub>c</sub>)-(hydroxyfunctional amide group) and wherein the (hydroxyfunctional amide group) is:

$$C$$
 $R_{4a}$ 

and  $R^{4a}$  is independently selected from the group consisting of OH,  $(C_1-C_6)$ alkoxy,  $(C_7-C_{14})$ alkaryloxy,  $(C_2-C_8)$ alkenyloxy,  $(C_7-C_{14})$  aralkyloxy,  $(C_7-C_{14})$ aralkenyloxy and aryloxy; and wherein  $R^{4b}$  is independently selected from the group consisting of H,  $(C_1-C_6)$ alkyl, arylalkyl, heteroaryl and aryl.

- 19. (Cancelled)
- 20. (Previously Presented) A compound selected from the group of:
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(hydroxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(methyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(methyl)-*N*-(methyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(hydroxy)-*N*-(methyl)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(ethyloxy)acetamide;

- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(2-propenyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(hydroxy)-N-(2-propyl)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(tert-butyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-[2-(methyl)propyloxy]acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(phenylmethyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(methyl)-*N*-(phenylmethyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-N-(phenyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(methyl)-*N*-(phenyloxy)acetamide;
- 2-[[3-(Aminooxoacetyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(cyclohexyl)-*N*-(hydroxy)acetamide; and
- 2-[[3-(2-Amino-2-oxoethyl)-2-ethyl-1-(phenylmethyl)-1H-indol-4-yl]oxy]-*N*-(hydroxy)acetamide.
  - 21. (Cancelled)
- 22. (Original) A pharmaceutical formulation comprising a indole compound as claimed in claim 1 together with a pharmaceutically acceptable carrier or diluent therefor.
  - 23. (Cancelled)
  - 24. (Cancelled)
- 25. (Previously Presented) A pharmaceutical formulation containing an effective amount of the compound of claim 1 useful for the treatment and/or amelioration of Inflammatory Diseases.
  - 26. (Cancelled)

Serial No. 10/629,992

27. (Cancelled)